INCLUDING INTERFERENCE

Ref #	Hits	Search Query	DBs	Default Operator	Plurals	Time Stamp
L1	2450	514/217.07 OR 514/249 OR 540/599 OR 544/349	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	OFF	2006/02/08 09:21
L2	647	L1 AND (ANTIPSYCHOTIC OR DOPAMINE OR SEROTONIN OR NOREPINEPHRINE OR EPINEPHRINE)	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	OFF	2006/02/08 09:22
L3	111	L2 AND (AZABICYCLO OR PYRIDOPYRAZINE OR PYRAZINOPYRIDINE OR ISOXAZOLE OR BENZOISOXAZOLE)	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	OFF	2006/02/08 09:22
L4	/1	L3 AND PYRIDYLOXYMETHYL	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	OFF	2006/02/08 09:23
L5	82	L3 AND (ANTAGONIST OR ANTAGONISTS)	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	OFF	2006/02/08 09:23
L6	√81	L5 NOT L4	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	OFF	2006/02/08 09:23

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chain nodes :

21-24 chain bonds: 6-11 21-22 21 ring bonds:

12-16 28-29 12-13 27-28 11-15 26-27 13-14 13-19 14-15 16-17 17-18 18-19 24-25 24-29 25-26

11-12 11-15 exact/norm bonds : 1-2 1-6 2-3 3-4 3-7 4-5 4-10 5-6 6-11 7-8 8-9 9-10 14-15 21-22 21-24 24-25 24-29 25-26 26-27 27-28 28-29 normalized bonds

13-14

12-16 13-19 16-17 17-18 18-19

isolated ring systems : containing 1 : 11 : 24 :

G1:C,N

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 19:Atom 19:Atom 19:Atom 21:CIASS 23:CIASS 24:Atom 25:Atom 26:Atom 27:Atom 28:Atom 29:Atom

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Structure attributes must be viewed using STN Express query preparation.

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13 ITERATIONS 100.0% PROCESSED SEARCH TIME: 00.00.01

13 ANSWERS

FULL FILE PROJECTIONS:

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44 TO 470 ONLINE BATCH PROJECTED ITERATIONS: PROJECTED ANSWERS:

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SINCE FILE ENTRY 166.94 cost in U.S. DOLLARS

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3 L4

D 1-3 IBIB ABS

Compositions containing atypical antipsychotics and azabicyclic compounds for treating CNS disorders Brodney, Michael A.; Howard, Harry R. Pfizer Inc, USA U.S. Pat. Appl. Publ., 21 pp. Patent CAPLUS COPYRIGHT 2006 ACS on STN 2005:698356 CAPLUS 143:179645 English FAMILY ACC. NUM. COUNT: PATENT INFORMATION: ANSWER 1 OF 3 PATENT ASSIGNEE(S) ACCESSION NUMBER: DOCUMENT NUMBER: TITLE: DOCUMENT TYPE: INVENTOR (S): SOURCE:

20050128 APPLICATION NO. A1 AM, AM, CCU, CCU, CCU, HR, LTT, LTT, TN, KE, KZ, KZ, SK, SK, TD, TD, PATENT NO.

of treating one or more CNS or other disorders, including concurrent treatment of disorders such as schizophrenia and depression. For example, eastbles for Parkinson's disease contained ziprasidone hydrochloride 200, benzisoxazole substituted azabicyclic compd 20, Methocel E3 222, lactose monohydrate 222, Aerosil 10, SLS 10 mg. method

CAPLUS COPYRIGHT 2006 ACS on STN
2004:780698 CAPLUS
141:296048
A preparation of pyrido[1,2-a]pyrazine derivatives, useful for the treatment of schizophrenia and ANSWER 2 OF 3 L5 ANSWER 2 OF 3 ACCESSION NUMBER: DOCUMENT NUMBER: TITLE:

depression
Tright, General: Brodney, Michael Aaron;
Wlodecki, Bishop
Pfizer Products Inc., USA
CODEN: PIXXDZ
CODEN: PIXXDZ INVENTOR (S):

PATENT ASSIGNEE (S)

Patent

English DOCUMENT TYPE:

APPLICATION NO. FAMILY ACC. NUM. COUNT: PATENT INFORMATION: PATENT NO.

BZ, CA, CH, FI, GB, GD, KR, KZ, LC, MZ, NA, NI, 20040223 ES, KP, MX, BR, BW, I EE, EG, I KE, KG, I WO 2004-IB499 8,5,5,8,8 BB, DZ, IS, BA, IN, AU, AZ, DE, DK, ID, IL, LV, MA, 20040923 A1 2 AM, AT, CU, CZ, HR, HU, LT, LU, 독유. 5. AE, AG, CN, CO, GE, GH, LK, LR, WO 2004081007

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TM, TM,
EW; EW; EW, KG,
ES, FT,
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CA 2518740
EP 1608648
R: AT, BE,
NL 1025710
NL 1025710 OTHER SOURCE(S):

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

The invefition relates to a preparation of pyrido[1,2-a]pyrazine derivs. of formula I [wherein: X is O, NH, or N(alkyl); Y is (CH2)0-1; Z is CHO, CO)-alkoxy, SO2-alkoxy, Me, CH2OH, etc.; R1 and R2 are independently selected from H, halogen, (cyclo)alkyl, or alkoxy, etc.], useful for treating CNS or other disorders, including concurrent treatment of disorders such as schizophrenia and depression. Thus, e.g., II was prepared via reaction of morpholine with (TR,985)-trans-methanesulfonic acid e-(2-benzo[d]isoxazol-3-yl-octahydropyrido[1,2-a]pyrazin-7-ylmethoxy)-pyridin-2-ylmethyl ester (preparation given). The prepared compds. were be antagonists and/or inverse agonists of human D2, human 5-HT1B, and human 5-HT2A receptors. For instance, preferred compound II exhibited Ki value of about 20 nM or less for at least two of the following receptors: D2, 5-HT1B, and 5-HT2A. determined to Æ

THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT REFERENCE COUNT

131:299376
Azabicyclic 5-HT1 receptor ligands, particularly 2-(benzo[d]isoxazol-3-y1)-7- $(phenoxymethy1) \, octahydropyrido \, [\, 1,2-a] \, pyrazine \, derivatives$ S COPYRIGHT 2006 ACS on STN 1999:672814 CAPLUS Bright, Gene Michael Pfizer Products Inc., USA PCT Int. Appl., 96 pp. CODEN: PIXXD2 15 ANSWER 3 OF 3 CAPLUS
ACCESSION NUMBER: 199
DOCUMENT NUMBER: 131
TITLE: Ara PATENT ASSIGNEE(S): INVENTOR(S): SOURCE:

English Patent FAMILY ACC. NUM. COUNT: PATENT INFORMATION: DOCUMENT TYPE:

Ξ 19990318 1, CZ, DE, JP, KE, MN, MW, TM, TR, RU, TJ, G, XX, XX, CN, IL, SL, KZ, CA, CH, HU, ID, LV, MD, SI, SK, BY, KG, APPLICATION NO. WO 1999-IB457 BY, HR, LU, SG, A.S. T. BG, GH, LLS, SD, ZW, 19991021 BA, BB, GD, GE, LK, LR, RO, RU, VN, YU, AZ, GB, LC, PT, KIND A1 AU, FI, KZ, PL, US, AT, ES, KR, NZ, UG, KP, KP, UA, AXXXXI PATENT NO. WO 9952907

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* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

MARPAT 131:299376

OTHER SOURCE(S): GI

The invention relates to compds. I [wherein R3, R4, Z = H, (fluoro)alkyl, (fluoro)alkoxy, (fluoro)alkoxy(fluoro)alkyl; W = alkoxymethyl, (fluoro)alkoxy(fluoro)alkyl; W = alkoxymethyl, (fluoro)alkoxy(fluoro)alkyl; W = alkoxymethyl, (fluoro)alkoxymethyl, or CHRNRIRZ where R1RZ = atoms to complete a heterocycle such as pyrrolidine; with provisos]. The compds. are agonists or antagonists of serotonin lA receptors, and/or antagonists of serotonin lD receptors, and are thus useful as psychotherapeutic agents. These compds. may be co-administered with 5-HT reuptake inhibitors, and are potentially useful for treating a wide variety of conditions. Approx. 40 synthetic examples are given. For instance, title compound II was prepared in 5 steps: (1) Mitsunobu etherification of starting material (7R,94s)-trans-III with Me 3-hydroxybenzoate (75%); (2) reduction of the Me ester to an alc. using LiAll4 (100%); (3) mesylation of the BCC protecting group (100%); and coupling with 3-chloro-5-fluoroescopic (36%). In assays against 5-HT receptors in vitro, all tested compds. I exhibited ICSO values of < 0.60 mM for 5-HTID receptors, and < 1.0 mM for 5-HTID REFERENCES AVAILABLE FOR THIS. DE 5-HTIA receptors. THERE ARE I CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

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ANSWER 3 OF 3 CAPLUS COPYRIGHT 2006 ACS on STN 247091-72-1P 247091-73-2P 247091-74-3P 247091-75-4P 247091-77-6P LS

NO PARTIMICAY METHY

GROUP IN ANY

(Reactant or reagent)
(intermediate; preparation of benzisoxazolyloctahydropyridopyrazine derivs, as 5-HT receptor ligands)
24/091-72-1 CAPLUS
Benzenemethanol, 3-[[(7R,9aS)-2-(5-fluoro-1,2-benzisoxazol-3-yl)octahydro-2H-pyrido[1,2-a]pyrazin-7-yl]methoxy]- (9CI) (CA INDEX NAME) Z Z

COLIPONING

Absolute stereochemistry.

Benzoic acid, 3-[[(7R,9aS)-2-(5-fluoro-1,2-benzisoxazol-3-y1)octahydro-2H-pyrido[1,2-a]pyrazin-7-y1]methoxy]-, methyl ester (9CI) (CA INDEX NAME) 247091-73-2 CAPLUS Z Z

Absolute stereochemistry.

247091-74-3 CAPLUS

Persenemethanol, 3-[[(7R,9aS]-2-(5-fluoro-1,2-benzisoxazol-3-yl)octahydroZH-pyridol1,2-a]pyrazin-7-yl]methoxy]-, methanesulfonate (ester) (9CI)
(CA INDEX NAME) S S

RN 247091-75-4 CAPLUS
CN Benzoic acid, 3-[[(75,9aS)-2-(1,2-benzisoxazol-3-y1)octahydro-2Hpyrido[1,2-a]pyrazin-7-y1]methoxy]-, methyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 247091-76-5 CAPLUS
CN Benzenemethanol, 3-[[(75,9aS)-2-(1,2-benzisoxazol-3-yl)octahydro-2H-pyrido[1,2-a]pyrazin-7-yl]methoxy]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 247091-77-6 CAPLUS
CN Benzcioc acid, 3-[[/7S,9aS]-2-(1,2-benzisoxazol-3-y])octahydro-2Hpyrido[1,2-a]pyrazin-7-yl]methoxy]-4-methyl-, methyl ester (9CI) (CA
INDEX NAME)

Absolute stereochemistry.

RN 247091-78-7 CAPLUS
CN Benzenemethanol, 3-[[(75,9aS)-2-(1,2-benzisoxazol-3-yl)octahydro-2H-pyrido[1,2-a]pyrazin-7-yl]methoxy]-4-methyl- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 247091-79-8 CAPLUS
CN Benzoic acid, 3-[[(7S,9aS)-2-(1,2-benzisoxazol-3-yl)octahydro-2Hpyrido[1,2-a]pyrazin-7-yl]methoxy]-5-methoxy-, methyl ester (9CI) (CA
INDEX NAME)

Absolute stereochemistry.

RN 247091-80-1 CAPLUS
CN Benzenemethanol, 3-[[(75,9aS)-2-(1,2-benzisoxazol-3-y1)octahydro-2H-pyrido[1,2-a]pyrazin-7-y1]methoxy]-5-methoxy- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 247091-81-2 CAPLUS
CN Benzoic acid, 5-[(175,9aS)-2-(1,2-benzisoxazol-3-yl)octahydro-2Hpyrido(1,2-a)pyrazin-7-yl)methoxy]-2-chloro-, methyl ester (9CI) (CA
INDEX NAME)

S S

247091-82-3 CAPLUS
Benzenemethanol, 5-[[(75,9aS]-2-(1,2-benzisoxazol-3-yl)octahydro-2H-pyrido[1,2-a]pyrazin-7-yl]methoxy]-2-chloro- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

₹ S

247091-83-4 CAPLUS
Benzoic acid, 4-[[(75,9aS)-2-(1,2-benzisoxazol-3-yl)octahydro-2Hpyrido[1,2-a]pyrazin-7-yl]methoxy]-, methyl ester (9Cl) (CA INDEX NAME)

Absolute stereochemistry.

₹ S

247091-84-5 CAPLUS
Benzenemethanol, 4-[[(7S,9aS)-2-(1,2-benzisoxazol-3-yl)octahydro-2H-pyrido[1,2-a]pyrazin-7-yl]methoxy]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

% S

247091-85-6 CAPLUS 2H-Pyrido[1,2-a]pyrazine, 2-(1,2-benzisoxazol-3-yl)-7-[(4-(chloromethyl)phenoxy]methyl]octahydro-, (7S,9aS)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

S 23

247091-88-9 CAPLUS
Benzonitrile, 4-[[(7S,9aS)-2-(1,2-benzisoxazol-3-yl)octahydro-2Hpyrido[1,2-a]pyrazin-7-yl]methoxy]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

S 25

247091-89-0 CAPLUS
Benzenemethanamine, 4-[[(7S,9aS)-2-(1,2-benzisoxazol-3-yl)octahydro-2H-pyrido[1,2-a]pyrazin-7-yl]methoxy]- (9CI) (CA INDEX NAME)

IT 247091-25-4P 247091-26-5P 247091-27-6P
247091-38-P 247091-32-BP 247091-33-4P
247091-31-2P 247091-32-BP 247091-33-4P
247091-31-2P 247091-33-3P 247091-33-4P
247091-34-5P 247091-33-9P 247091-39-0P
247091-40-3P 247091-44-7P 247091-42-5P
247091-40-3P 247091-44-7P 247091-42-PP
247091-50-5P 247091-44-7P 247091-42-PP
247091-50-5P 247091-51-6P 247091-52-7P
247091-50-5P 247091-47-P 247091-47-P
247091-50-5P 247091-47-P
247091-50-5P 247091-47-P
247091-47-H

derivs.

as 5-HT1 receptor ligands)
247091-25-4 CRPLUS
Benzenemethanamine, 3-[[(7S,9aS)-2-(1,2-benzisoxazol-3-yl)octahydro-2Hpyrido[1,2-a]pyrazin-7-yl]methoxy]-N-cyclopropyl- (9CI) (CA INDEX NAME) S S

Absolute stereochemistry.

S &

247091-26-5 CAPLUS
24-Pyridol[1,2-a]pyrazine, 2-(1,2-benzisoxazol-3-yl)octahydro-7-[[3-[[(2S)-2-(methoxymethyl)-1-pyrrolidinyl]methyl]phenoxy]methyl]-, (7S,9aS)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

247091-27-6 CAPLUS
2H-Pyrido[1,2-a]pyrazine, 2-(1,2-benzisoxazol-3-y1)-7-[[3-[(4-ethy1-1-pperaziny1)methy1]phenoxy]methy1]octahydro-, (75,9aS)- (9CI) (CA INDEX NAME) S S

Absolute stereochemistry.

247091-28-7 CAPLUS 2H-Pyrido[1,2-a]pyrazine, 2-(5-fluoro-1,2-benzisoxazol-3-yl)octahydro-7-[[3-(1-pyrrolidinylmethyl)phenoxy]methyl]-, (7R,9aS)- (9CI) (CA INDEX NAME) S S

Absolute stereochemistry.

247091-29-8 CAPLUS
3-Azetidinol, 1-[[3-[[/7k,9aS]-2-(5-fluoro-1,2-benzisoxazol-3-yl)octahydro2H-pyrido[1,2-a]pyrazin-7-yl]methoxylphenyl]methyl]- (9CI) (CA INDEX
NAME) Z Z

S S

247091-30-1 CAPLUS
24-9yrido[1,2-a]pyrazine, 2-(5-fluoro-1,2-benzisoxazol-3-yl)octahydro-7[[3-(4-morpholinylmethyl)phenoxy]methyl]-, (7R,9aS)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

S Z

247091-31-2 CAPLUS
2H-Pyrido[1,2-a]pyrazine, 2-(4-fluoro-1,2-benzisoxazol-3-yl)octahydro-7[[3-(1-pyrrolidinylmethyl)phenoxy]methyl]-, (7R,9aS)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

Z Z

247091-32-3 CAPLUS
24-Pyrido[1,2-a]pyrazine, 2-(1,2-benzisoxazol-3-yl)octahydro-7-[[3-(1-pyrrolidinylmethyl)phenoxy]methyl]-, (7S,9aS)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

S S

247091-33-4 CAPLUS
A+Pyrrol.idinediol, 1-[[3-[[178,9as]-2-(1,2-benzisoxazol-3-y1)octahydro-2H-pyrido[1,2-a]pyrazin-7-y1]methoxy]phenyl]methyl]-, (3S,4S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

247091-34-5 CAPLUS 2H-Pyrido[1,2-a]pyrazine, 2-(1,2-benzisoxazol-3-yl)octahydro-7-[[2-methyl-5-(1-pyrrolidinylmethyl)phenoxy]methyl]-, (7S,9aS)- (9CI) (CA INDEX NAME) Z Z

Absolute stereochemistry.

247091-35-6 CAPLUS
2H-Pyrido[1,2-a]pyrazine, 2-(1,2-benzisoxazol-3-y1)octahydro-7-[[3-methoxy-5-(1-pyrrolidinylmethy1)phenoxy]methy1]-, (78,9a8)- (9CI) (CA INDEX NAME) Z Z

RN 247091-36-7 CAPLUS
CN 2H-Pyrido(1,2-a)pyrazine, 2-(1,2-benzisoxazol-3-y1)-7-[(4-chloro-3-(1-pyrrolidinylmethyl)phenoxy]methyl]octahydro-, (75,9as)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 247091-37-8 CAPLUS
CN 2H-Pyrido[1,2-a]pyrazine, 2-(1,2-benzisoxazol-3-yl)octahydro-7-[[4-(1-pyrrolidinylmethyl)phenoxy]methyl]-, (75,9aS)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 247091-38-9 CAPLUS
CN 2H-Pyrido[1,2-a]pyrazine, 7-[[3-(1-azetidinylmethyl)phenoxy]methyl]-2-(1,2-benzisoxazoi-3-yl)octahydro-, (78,9aS)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 247091-39-0 CAPLUS
CN Benzenemethanamine, 3-[[(75,9as)-2-(1,2-benzisoxazol-3-yl)octahydro-2H-pyrido[1,2-a]pyrazin-7-yl]methoxy]-N-(cyclopropylmethyl)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 247091-40-3 CAPLUS
CN Benzenemethanamine, 3-[[(78,9aS)-2-(1,2-benzisoxazol-3-y1)octahydro-2H-pyrido[1,2-a]pyrazin-7-y1]methoxy]-N-cyclohexy1- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 247091-41-4 CAPLUS CN 3-Pyrrolidinol, 1-[(3-[(75,9as)-2-(1,2-benzisoxazol-3-yl)octahydro-2H-pyrido[1,2-a]pyrazin-7-yl]methoxy]phenyl]methyl]-, (3R)- (9CI) (CA INDEX NAME)

Z Z

247091-42-5 CAPLUS
ZH-PYLOIG(1,2-3) pyrazine, 2-(1,2-benzisoxazol-3-yl)-7-[(5-([(25,55)-2,5-dimethyl-1-pyrrolidinyl]methyl]-2-methylphenoxy]methyl]octahydro-, (75,9as)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

247091-43-6 CAPLUS
2H-PY1040(1,2-a) pyrazine, 2-(1,2-benzisoxazol-3-yl)-7-[(5-([(2R,5R)-2,5-dimethyl-1-pyrrolidinyl]methyl)-2-methylphenoxy]methyl]octahydro-, (7S,9aS)- (9CI) (CA INDEX NAME) S S

Absolute stereochemistry.

247091-44-7 CAPLUS
3-Pyrrolidinol, 1-[(3-[(78,9aS)-2-(1,2-benzisoxazol-3-yl)octahydro-2H-pyrido[1,2-ajpyrazin-7-yl]methoxy]phenyl]methyl]-, (3S)- (9CI) (CA INDEX NAME) Z Z

Absolute stereochemistry.

247091-45-8 CAPLUS

Benzenemethanamine, 3-[[/7S,9aS]-2-(1,2-benzisoxazol-3-yl)octahydro-2Hpyrido[1,2-a]pyrazin-7-yl]methoxy]-N-(2-methylpropyl)- (9CI) (CA INDEX NAME) G &

Absolute stereochemistry.

247091-47-0 CAPLUS
2H-Pyrido[1,2-a]pyrazine, 2-(1,2-benzisoxazol-3-yl)octahydro-7-[{2-(1-pyrrolidinylmethyl)phenoxy]methyl]-, (7S,9aS)- (9CI) (CA INDEX NAME) Z Z

Absolute stereochemistry.

247091-48-1 CAPLUS
24-Pyrido[1,2-a]pyrazine, 2-(1,2-benzisoxazol-3-yl)octahydro-7-[[4-(4-morpholinylmethyl)phenoxylmethyl]-, (7S,9aS)- (9CI) (CA INDEX NAME) S S

RN 247091-49-2 CAPLUS
CN 2H-Pyrido[1,2-a]pyrazine, 2-(7-fluoro-1,2-benzisoxazol-3-yl)octahydro-7-[[3-(1-pyrrolidinylmethyl)phenoxy]methyl]-, (7R,9aS)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 247091-50-5 CAPLUS
CN 2H-Pyrido[1,2-a]pyrazine, 2-(6-fluoro-1,2-benzisoxazol-3-yl)octahydro-7[{3-(1-pyrrolidinylmethyl)phenoxylmethyl]-, (7R,9aS)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 247091-51-6 CAPLUS
CN 2H-Pyrido[1,2-a]pyrazine, 2-(6,7-difluoro-1,2-benzisoxazol-3-y1)octahydroT-[[3-(1-pyrrolidinylmethyl)phenoxy]methyl]-, (7R,9aS)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 247091-52-7 CAPLUS
CN 3-Azabicyclo[3.2.2]nonane, 3-[[3-[[(7R,9aS)-2-(5-fluoro-1,2-benzisoxazol-3-y])octahydro-2H-pyrido[1,2-a]pyrazin-7-yl]methoxylphenyl]methyl]- (9CI)
(CA INDEX NAME)

Absolute stereochemistry.

RN 247091-53-8 CAPLUS
CN 2H-Pyrido[1,2-a]pyrazine, 2-(5-fluoro-1,2-benzisoxazol-3-yl)octahydro-7[[3-[(octahydro-2H-isoindol-2-yl)methyl]phenoxy]methyl]-, (7R,9aS)- (9CI)
(CA INDEX NAME)

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